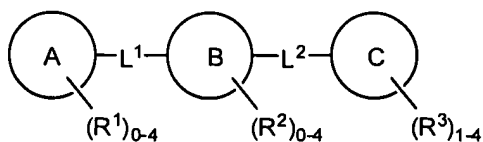


Listing of the Claims:

1. (Currently Amended) A compound for modulating c-Kit activity according to Formula I,



I

or a pharmaceutically acceptable salt, thereof, wherein,

ring A is:

wherein each Y is independently either =C(H)- or =N-; and Z is selected from -O-, -S-, and

-N(R⁷)-;

each R¹ is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two adjacent of R¹, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R¹⁰;

L¹ is a single bond;

ring B is phenyl ;

each R² is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴, -N(R⁴)R⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two adjacent of R², together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R¹⁵;

L² is selected from -N(H)N(H)C(=O)N(H)-, -CH₂N(H)C(=O)N(H)-, -CH₂OC(=O)N(H)-, and -XCH₂C(=O)N(H)-; wherein X is selected from -O-, -S(O)₀₋₂-, and -N(R⁷)-; and any C-H of L² is optionally C-R²⁰;

ring C is phenyl ;

each R³ is independently selected from halogen, trihalomethyl, -CN,

-NO₂, -OR⁴, -N(R⁴)R⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴,

$-\text{C}(=\text{NR}^5)\text{R}^4$, $-\text{N}(\text{R}^4)\text{SO}_2\text{R}^4$, $-\text{N}(\text{R}^4)\text{C}(\text{O})\text{R}^4$, $-\text{NCO}_2\text{R}^4$, $-\text{C}(=\text{O})\text{R}^4$, optionally substituted alkoxy, optionally substituted C_{1-6} alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6} alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6} alkyl; provided R^3 is not a cyclic sulfonamide attached to ring C via the nitrogen of said cyclic sulfonamide,

wherein there exists at least one of R^3 that is halogen or trihalomethyl;

R^4 is selected from -H, optionally substituted C_{1-6} alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6} alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6} alkyl;

two of R^4 , when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R^5 is selected from -H, -CN, -NO₂, -OR⁴, -S(O)₀₋₂R⁴, -CO₂R⁴, optionally substituted C_{1-6} alkyl, optionally substituted C_{1-6} alkenyl, and optionally substituted C_{1-6} alkynyl;

R^7 is selected from -H, optionally substituted C_{1-6} alkyl, -SO₂N(R⁴)R⁴, -CO₂R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted aryl C_{1-6} alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6} alkyl; and

each of R^{10} , each of R^{15} , each of R^{20} , and each of R^{25} is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴, -N(R⁴)R⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -CO₂R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C_{1-6} alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6} alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6} alkyl;

provided:

~~when both ring B and ring C are phenyl;~~

- a) ~~and the compound comprises ring B-CH₂N(H)C(=O)N(H)-ring C, then L¹ must be a single bond; R² can not comprise a group of the formula -O(CH₂)_{2,4}-N-piperazine that is *ortho* to L²; and ring A cannot be a 5-methyl-[1,2,4]-oxadiazol-3-yl radical, a 4H-[1,2,4]-oxadiazol-5-one-3-yl radical, nor a 4'-[2,2';6',2'']terpyridinyl radical;~~
- b) ~~and L¹ is single bond, then L² cannot comprise N(H)C(=O)C(=O)N(H) nor N(H)C(=Q)C(H)CNC(=O) (where Q is S or O);~~
- c) ~~and L¹ is other than single bond, then A cannot be quinolin-2-yl-L¹, quinolin-3-yl-L¹, or quinolin-4-yl-L¹;~~

~~;~~ and

the compound is not one of: N-naphthalen-1-yl-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(phenyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,6-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-ethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,6-diethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(ethyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(dimethylamino)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{{[3-(1H-tetrazol-1-

~~yl)phenyl]oxy}acetamide, N-(4-chloro-3-methylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-fluorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoromethyl)phenyl]acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl)phenyl]acetamide, methyl-4-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoate, ethyl-4-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoate, 3-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoic acid, N-[3-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy}acetamide, acetamide, acetamide, or N-(4-chlorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-aminophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, and N-(4-acetylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide.~~

2-12. (cancelled)

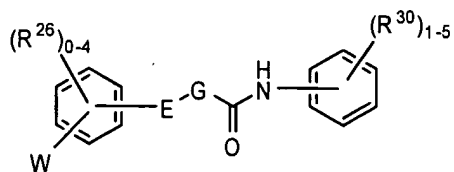
13. (Currently Amended) The compound according to ~~claim 10~~, claim 1, wherein there exists at least one of R³ that is trifluoromethyl.

14. (original) The compound according to claim 13, wherein ring C is a phenyl comprising a trifluoromethyl radical *meta*- to L².

15. (Previously Presented) The compound according to claim 1, wherein each of R³ is independently selected from halogen, trihalomethyl, -OR⁴, -C(=O)R⁴, and optionally substituted C₁₋₆alkyl.

16. (Previously Presented) A compound for modulating c-Kit activity according to the

following Formula:



or a pharmaceutically acceptable salt, thereof, wherein,

W is selected from the following:

each of R^{27} independently selected from halogen, trihalomethyl, -CN, -NO₂, -OR⁵⁵, -S(O)₀₋₂R⁵⁵, -SO₂N(R⁵⁵)R⁵⁵, -C(=O)N(R⁵⁵)R⁵⁵, -C(=NR⁵⁰)N(R⁵⁵)R⁵⁵, -C(=NR⁵⁰)R⁵⁵, -N(R⁵⁵)SO₂R⁵⁵, -N(R⁵⁵)C(O)R⁵⁵, -NCO₂R⁵⁵, -C(=O)R⁵⁵, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

each Y is independently either =C(H)- or =N-;

Z is selected from -O-, -S(O)₀₋₂-, and -N(R⁷)-

E and G are each independently selected from -O-, -S(O)₀₋₂-, -C(R³¹)R³²-, and -N(R³³)-

J₁ and J₂ are each independently =C(H)- or =N-;

R²⁶ is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -CO₂R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -N(R⁴⁰)SO₂R⁴⁰, -N(R⁴⁰)C(O)R⁴⁰, -NCO₂R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

R³⁰ is independently selected from halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -N(R⁴⁰)SO₂R⁴⁰, -N(R⁴⁰)C(O)R⁴⁰, -NCO₂R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl, wherein there exists at least one of R³⁰ that is trihalomethyl;

R³¹ and R³² are each independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -CO₂R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -N(R⁴⁰)SO₂R⁴⁰, -N(R⁴⁰)C(O)R⁴⁰, -NCO₂R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

R³³ is selected from -H, optionally substituted lower alkyl, -SO₂N(R⁴⁰)R⁴⁰, -CO₂R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

R⁴⁰ is selected from -H, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two of R⁴⁰, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R⁵⁰ is selected from -H, -CN, -NO₂, -OR⁴⁰, -S(O)₀₋₂R⁴⁰, -CO₂R⁴⁰, optionally substituted C₁₋₆alkyl, optionally substituted C₁₋₆alkenyl, and optionally substituted C₁₋₆alkynyl;

R⁵⁵ is selected from -H, optionally substituted C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl; and

two of R⁵⁵, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P.

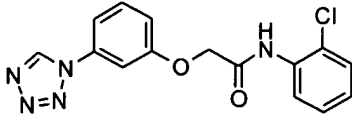
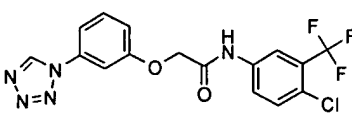
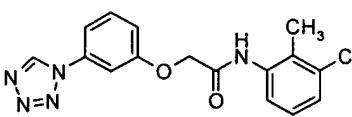
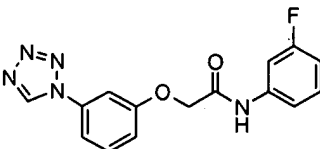
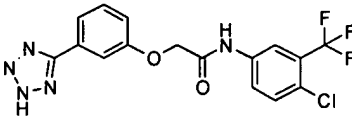
17. (cancelled)

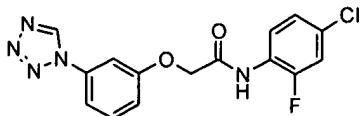
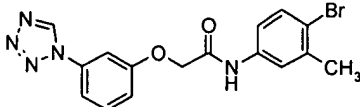
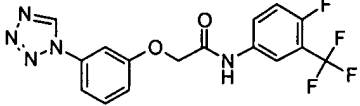
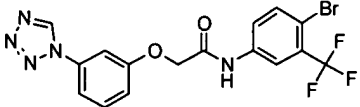
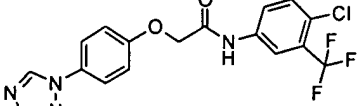
18. (previously presented) The compound according to claim 16, wherein R³⁰ is selected from halogen, trihalomethyl, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl, wherein there exists at least one of R³⁰ that is trifluoromethyl.

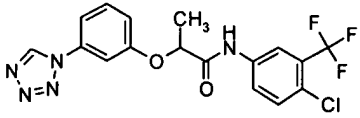
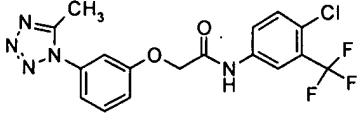
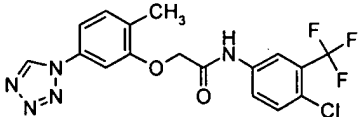
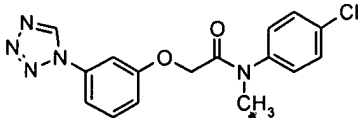
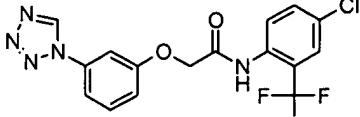
19. (cancelled)
20. (cancelled)
21. (cancelled)
22. (cancelled)
23. (withdrawn from consideration, currently amended) The compound according to ~~claim 22~~, claim 16, wherein E is selected from -O-, -S(O)₀₋₂-, and -NH-; and G is -CH₂-.
24. (withdrawn from consideration, currently amended) The compound according to ~~claim 22~~, claim 16, wherein E is either -CH₂- or -NH-; and G is selected from -O-, -S-, and -NH-.
25. (cancelled)
26. (cancelled)
27. (currently amended)

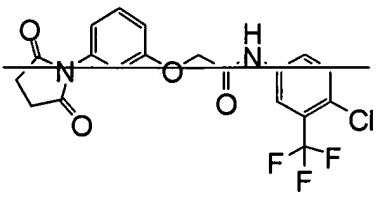
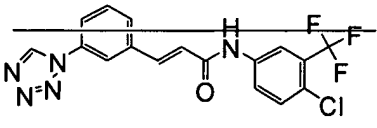
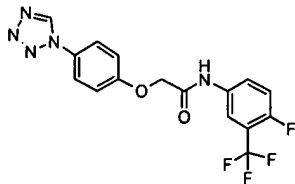
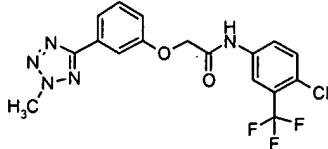
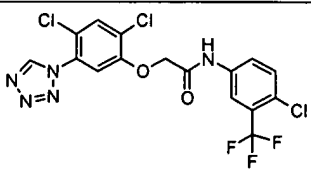
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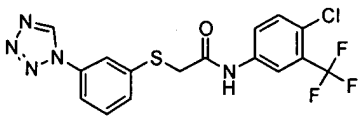
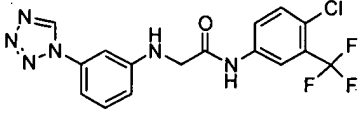
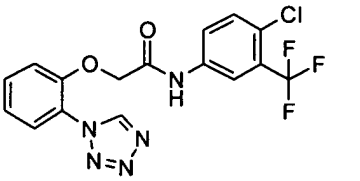
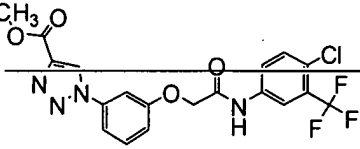
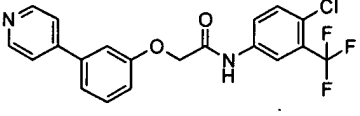
Entry	Name	Structure
1	N-[5-chloro-2-(methoxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Entry	Name	Structure
4	N-(2-chlorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
5	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
7	N-(3-chloro-2-methylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
8	N-(3-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
9	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(2H-tetrazol-5-yl)phenyl}oxy}acetamide	

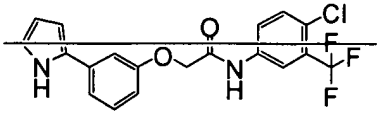
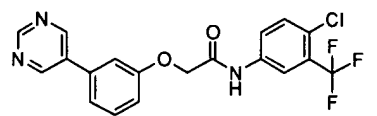
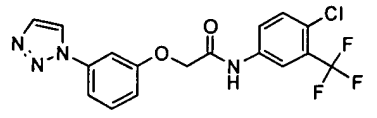
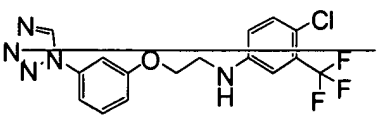
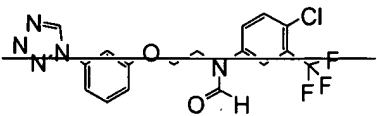
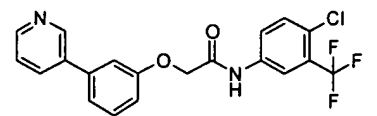
Entry	Name	Structure
10	N-(4-chloro-2-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
11	N-(4-bromo-3-methylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
12		
13	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
14	N-[4-bromo-3-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
15	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{4-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	

Entry	Name	Structure
16	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} propanamide	
17	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(5-methyl-1H-tetrazol-1-yl)phenyl]oxy} acetamide	
18	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-methyl-5-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
19	N-(4-chlorophenyl)-N-methyl-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
20	N-[4-chloro-2-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	

Entry	Name	Structure
21	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(2,5-dioxopyrrolidin-1-yl)phenyl]oxy}acetamide	
22	(2E)-N-[4-chloro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]prop-2-enamide	
23	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2- {[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
24	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(2-methyl-2H-tetrazol-5-yl)phenyl]oxy}acetamide	
25	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2,4-dichloro-5-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Entry	Name	Structure
26	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]thio}acetamide	
27	N-[4-chloro-3-(trifluoromethyl)phenyl]- N~2~-[3-(1H-tetrazol-1- yl)phenyl]glycinamide	
28	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	
29	methyl 1-{3-[(2-{[4-chloro-3- (trifluoromethyl)phenyl]amino}-2- oxoethyl)oxy]phenyl}-1H-1,2,3-triazole-4- carboxylate	
37	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-pyridin-4-ylphenyl)oxy]acetamide	

Entry	Name	Structure
38	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-methyl-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
48	5-chloro-2-(((3-(1H-tetrazol-1-yl)phenyl)oxy}acetyl)amino}benzamide	
49	N-[5-chloro-2,4-bis(methoxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
52	N-[2-(methoxy)-5-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
57	1,1-dimethylethyl 2-{3-[(2-[[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-pyrrole-1-carboxylate	

Entry	Name	Structure
58	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-pyrrol-2-yl)phenyl]oxy}acetamide	
59	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-pyrimidin-5-yl)phenyl]oxy}acetamide	
60	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-1,2,3-triazol-1-yl)phenyl]oxy} acetamide	
61	4-chloro-N-(2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}ethyl)-3-(trifluoromethyl)aniline	
62	N-[4-chloro-3-(trifluoromethyl)phenyl]-N-(2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}ethyl)formamide	
63	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-pyridin-3-yl)phenyl]oxy}acetamide	

Entry	Name	Structure
64	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-furan-3-ylphenyl)oxy]acetamide	
65	(2E)-N-[4-fluoro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]prop-2-enamide	
66	N-[4-fluoro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]propanamide	
68	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(3,5-dimethylisoxazol-4-yl)phenyl]oxy}acetamide	
69	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-quinolin-7-ylphenyl)oxy]acetamide	
70	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-furan-2-ylphenyl)oxy]acetamide	

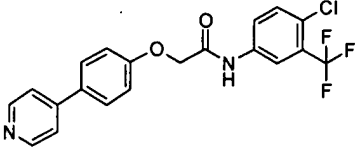
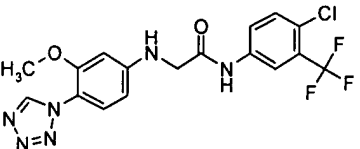
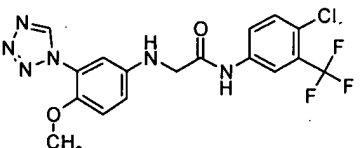
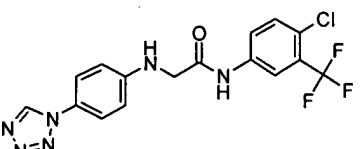
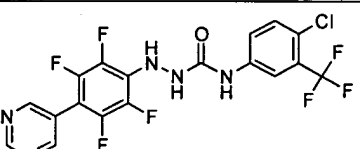
Entry	Name	Structure
71	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
72	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-dibenzo[b,d]furan-4-ylphenyl)oxy]acetamide	
73	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	
75	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-tetrazol-1-yl)phenyl]methyl}urea	
76	N-[4-chloro-3-(trifluoromethyl)phenyl]-N-methyl-2-[(3-(1H-tetrazol-1-yl)phenyl)oxy]acetamide	

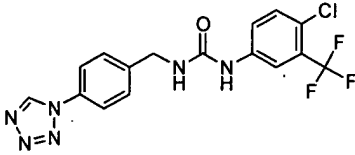
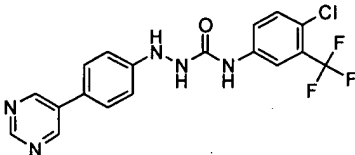
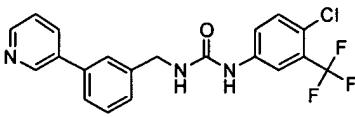
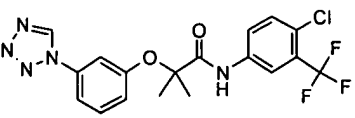
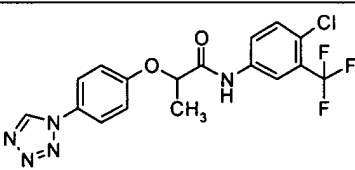
Entry	Name	Structure
77	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
78	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[[3-(pyridin-2-ylamino)phenyl]oxy]acetamide	
79	N-[2-fluoro-5-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
80	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	
81	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea	

Entry	Name	Structure
82	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyrimidin-5-ylphenyl)methyl]urea	
83	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
84	[3-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
85	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	
86	N-2-[4-chloro-3-(trifluoromethyl)phenyl]-N-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	

Entry	Name	Structure
87	2-{{4-chloro-3-(trifluoromethyl)phenyl}oxy}-N-[3-(1H-tetrazol-1-yl)phenyl]acetamide	
88	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-methyl-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
89	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[4-(1H-1,2,3-triazol-1-yl)phenyl]oxy}acetamide	
90	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
91	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[2-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

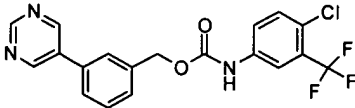
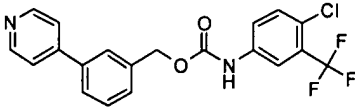
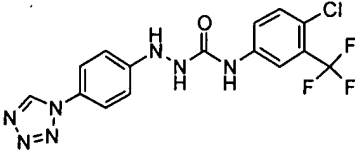
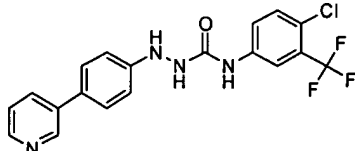
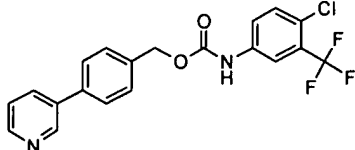
Entry	Name	Structure
92	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-3-(1H-tetrazol-1-yl)benzenesulfonamide	
93	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-N-methyl-3-(1H-tetrazol-1-yl)benzenesulfonamide	
94	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	
95	2-({4-[2,4-bis(methoxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-fluoro-3-(trifluoromethyl)phenyl]acetamide	
96	2-({4-[2,4-bis(methoxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-chloro-3-(trifluoromethyl)phenyl]acetamide	

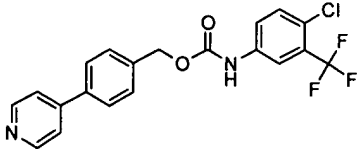
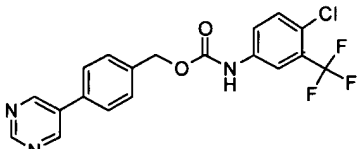
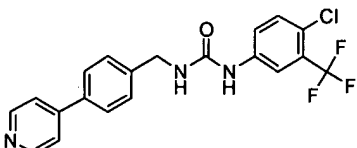
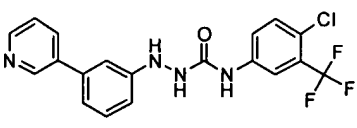
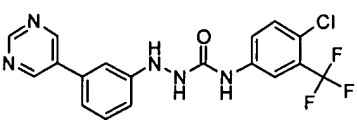
Entry	Name	Structure
97	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-4-ylphenyl)oxy]acetamide	
98	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(methoxy)-4-(1H-tetrazol-1-yl)phenyl]glycinamide	
99	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(methoxy)-3-(1H-tetrazol-1-yl)phenyl]glycinamide	
100	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(1H-tetrazol-1-yl)phenyl]glycinamide	
101	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(2,3,5,6-tetrafluoro-4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	

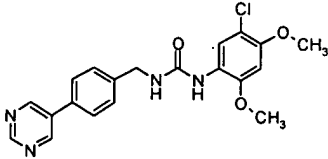
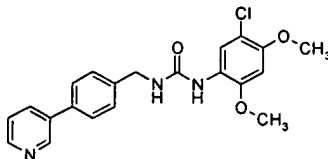
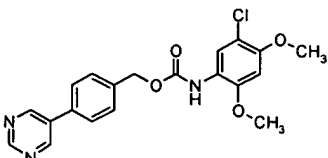
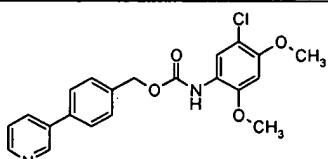
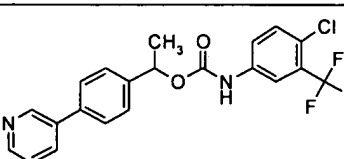
Entry	Name	Structure
102	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-tetrazol-1-yl)phenyl]methyl}urea	
103	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
104	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyridin-3-ylphenyl)methyl]urea	
105	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
106	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	

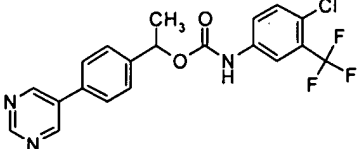
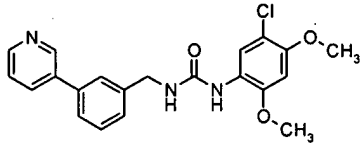
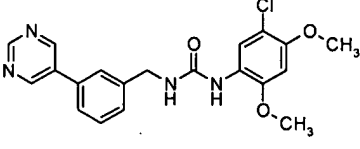
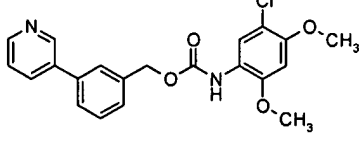
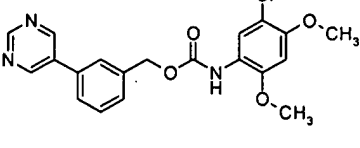
Entry	Name	Structure
107	N-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
108	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[2-(methyloxy)pyrimidin-5-yl]phenyl}methyl)urea	
109	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
110	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[2-(methyloxy)pyrimidin-5-yl]phenyl}methyl)urea	
111	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	

Entry	Name	Structure
112	1,1-dimethylethyl 2-{4-[(2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-indole-1-carboxylate	
113	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-4-(1H-tetrazol-1-yl)benzenesulfonamide	
114	N-[4-chloro-3-(trifluoromethyl)phenyl]-N-2-[3-(2H-tetrazol-5-yl)phenyl]glycinamide	
115	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[2,6-difluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
116	(3-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

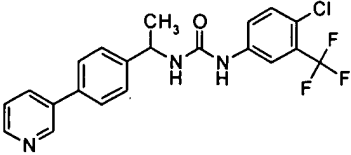
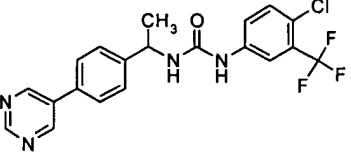
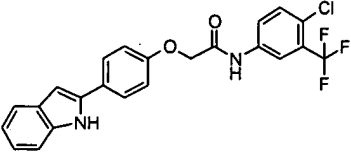
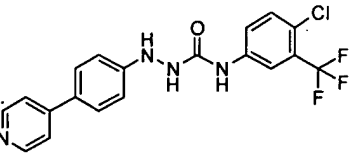
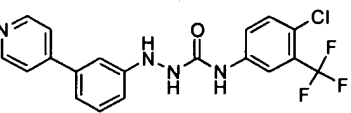
Entry	Name	Structure
117	(3-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
118	(3-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
119	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[4-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
120	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-3-ylphenyl)hydrazinecarboxamide	
121	(4-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Entry	Name	Structure
122	(4-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
123	(4-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
124	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-4-ylphenyl)methyl]urea	
125	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-3-ylphenyl)hydrazinecarboxamide	
126	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide	

Entry	Name	Structure
127	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(4-pyrimidin-5-yl)phenyl)methyl]urea	
128	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(4-pyridin-3-yl)phenyl)methyl]urea	
129	(4-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	
130	(4-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	
131	1-(4-pyridin-3-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Entry	Name	Structure
132	1-(4-pyrimidin-5-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
133	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(3-pyridin-3-ylphenyl)methyl]urea	
134	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea	
135	(3-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	
136	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	

Entry	Name	Structure
137	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
138	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
139	N-[[3-(6-aminopyridin-3-yl)phenyl]methyl]-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
140	N-[[4-(6-aminopyridin-3-yl)phenyl]methyl]-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
141	N-[[3-(2-aminopyrimidin-5-yl)phenyl]methyl]-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
142	N-[[4-(2-aminopyrimidin-5-yl)phenyl]methyl]-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Entry	Name	Structure
143	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyridin-3-ylphenyl)ethyl]urea	
144	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyrimidin-5-ylphenyl)ethyl]urea	
145	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-indol-2-yl)phenyl]oxy}acetamide	
147	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-4-ylphenyl)hydrazinecarboxamide	
148	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-4-ylphenyl)hydrazinecarboxamide	

Entry	Name	Structure
149	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyridin-4-ylphenyl)methyl]urea	
150	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-quinoxalin-6-ylphenyl)methyl]urea	
151	methyl 3-amino-6-(3-(((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methylphenyl)pyrazine-2-carboxylate	
152	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-quinoxalin-6-ylphenyl)methyl]urea	
153	N-([3-(2-amino-5-methylpyridin-3-yl)phenyl]methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

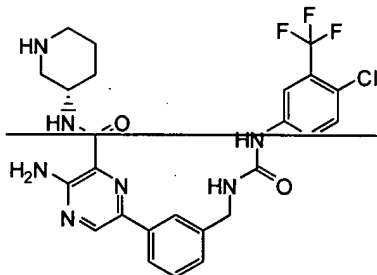
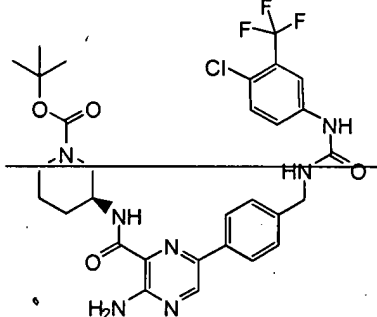
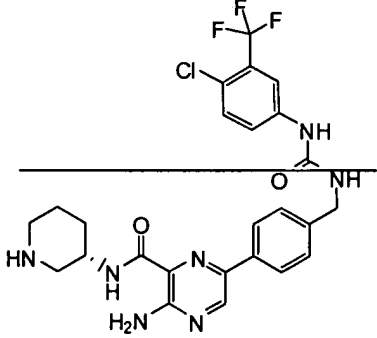
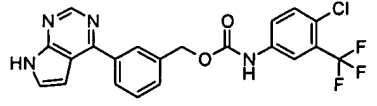
Entry	Name	Structure
154	methyl 3-amino-6-(4-(((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)pyrazine-2-carboxylate	
155	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-chloro-4-(methyloxy)phenyl]carbamate	
156	N-[3-chloro-4-(methyloxy)phenyl]-N'-{[3-(1H-tetrazol-1-yl)phenyl]methyl}urea	
157	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[4-(5-hydroxy-1H-tetrazol-1-yl)phenyl]oxy}acetamide	
158	N-{{[3-(2-amino-5-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
159	N-{{[4-(2-amino-5-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

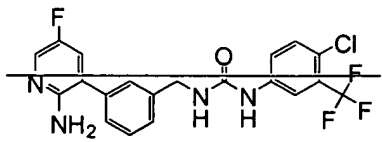
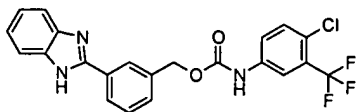
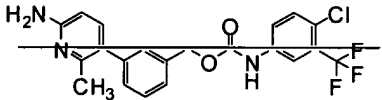
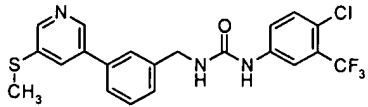
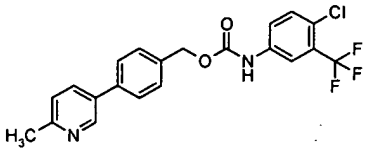
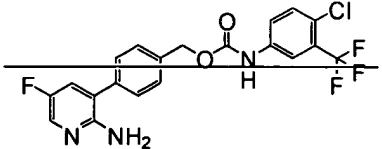
Entry	Name	Structure
160	N-{[3-(6-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
161	N-{[4-(6-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
162	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(pyrimidin-2-yloxy)phenyl]methyl}urea	
163	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-3-(1H-tetrazol-1-yl)benzamide	
164	3-amino-6-(3-{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)-N-[2-(dimethylamino)ethyl]pyrazine-2-carboxamide	
165	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-fluoropyridin-3-yl)phenyl]methyl}urea	

Entry	Name	Structure
166	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({3-[2-(methyloxy)pyridin-3- yl]phenyl}methyl)urea	
167	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(6-fluoropyridin-3- yl)phenyl]methyl}urea	
168	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({4-[2-(methyloxy)pyridin-3- yl]phenyl}methyl)urea	
169	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(6-methylpyridin-3- yl)phenyl]methyl}urea	
170	N-{{4-(2-amino-5-fluoropyridin-3- yl)phenyl}methyl}-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea	

Entry	Name	Structure
171	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[3-(6-methylpyridin-3-yl)phenyl]methyl}urea	
172	N- {[4-(2-aminopyridin-3-yl)phenyl]methyl} N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
173	N- {[3-(2-aminopyridin-3-yl)phenyl]methyl} N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
174	[3-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
175	[3-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
176	[3-(2-aminopyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Entry	Name	Structure
177	(3-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
178	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({[3-(6-(hydroxymethyl)pyridin-3-yl)phenyl]methyl})urea	
179	N-{[3-(6-acetylpyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
180	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-cyanopyridin-3-yl)phenyl]methyl}urea	
181	1,1-dimethylethyl (3S)-3-({[3-amino-6-(3-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)pyrazin-2-yl]carbonyl)amino)piperidine-1-carboxylate	

Entry	Name	Structure
182	3-amino-6-(3-(((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)-N-((3S)-piperidin-3-yl)pyrazine-2-carboxamide	
183	1,1-dimethylethyl (3S)-3-(((3-amino-6-(4-(((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)pyrazin-2-yl)carbonyl)amino)piperidine-1-carboxylate	
184	3-amino-6-(4-(((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)-N-((3S)-piperidin-3-yl)pyrazine-2-carboxamide	
185	[3-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Entry	Name	Structure
186	N-[[3-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl]-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
188	[3-(1H-benzimidazol-2-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
189	[3-(6-amino-2-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
190	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-((3-[5-(methylthio)pyridin-3-yl]phenyl)methyl)urea	
191	[4-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
192	[4-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

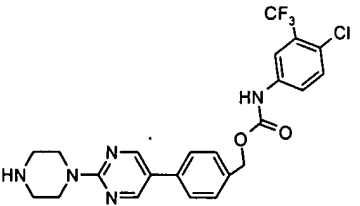
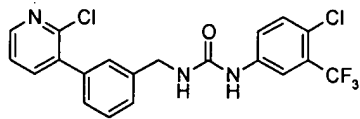
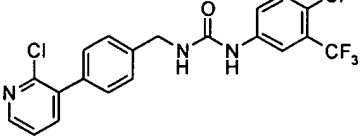
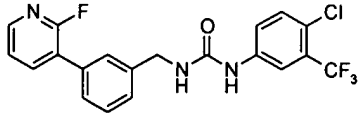
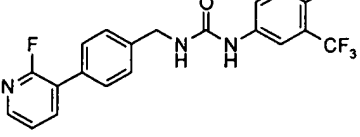
Entry	Name	Structure
193	[4-(2-aminopyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
194	(4-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
195	[4-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
196	[4-(6-amino-2-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
199	(3-pyridin-3-ylphenyl)methyl (3,5-dimethylphenyl)carbamate	
200	(3-pyridin-3-ylphenyl)methyl [5-chloro-2-(methyloxy)phenyl]carbamate	

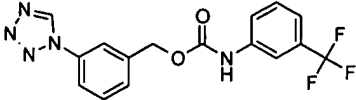
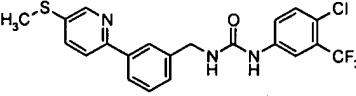
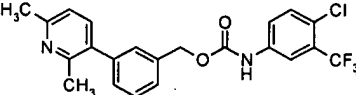
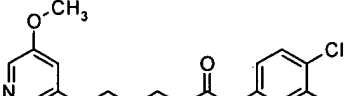
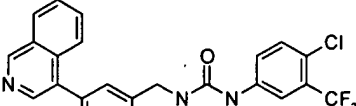
Entry	Name	Structure
201	[4-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
202	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2-(methoxy)phenyl]carbamate	
203	(4-pyrimidin-5-ylphenyl)methyl (3,4-dimethylphenyl)carbamate	
204	(3-pyridin-3-ylphenyl)methyl (3,4-dimethylphenyl)carbamate	
205	1,1-dimethylethyl 3-(((3-amino-6-(3-(((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)pyrazin-2-yl)carbonyl)amino)piperidine-1-carboxylate	

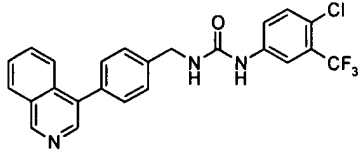
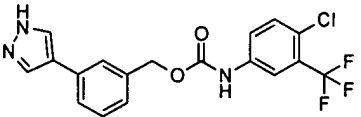
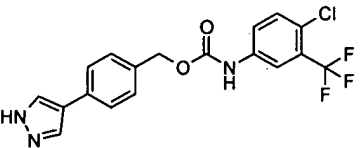
Entry	Name	Structure
206	1,1-dimethylethyl 3-(((3-amino-6-(4-(((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)pyrazin-2-yl)carbonyl)amino)piperidine-1-carboxylate	
207	3-amino-6-(3-(((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)-N-piperidin-3-ylpyrazine-2-carboxamide	
208	3-amino-6-(4-(((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)-N-piperidin-3-ylpyrazine-2-carboxamide	

Entry	Name	Structure
209	1,1 dimethylethyl 4-{{3-amino-6-(3- (((4-chloro-3- (trifluoromethyl)phenyl)amino)carbonyl)amino)methyl}phenyl)pyrazin-2-yl}carbonyl}piperazine-1-carboxylate	
210	1,1 dimethylethyl 4-{{3-amino-6-(4- (((4-chloro-3- (trifluoromethyl)phenyl)amino)carbonyl)amino)methyl}phenyl)pyrazin-2-yl}carbonyl}piperazine-1-carboxylate	
211	N-({3-[5-amino-6-(piperazin-1-ylcarbonyl)pyrazin-2-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Entry	Name	Structure
212	N-({4-[5-amino-6-(piperazin-1-ylcarbonyl)pyrazin-2-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
213	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-pyrazol-4-yl)phenyl]methyl}urea	
214	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-pyrazol-4-yl)phenyl]methyl}urea	
215	[3-(2-piperazin-1-yl)pyrimidin-5-yl]phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Entry	Name	Structure
216	[4-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
217	N-{[3-(2-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
218	N-{[4-(2-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
219	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(2-fluoropyridin-3-yl)phenyl]methyl}urea	
220	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(2-fluoropyridin-3-yl)phenyl]methyl}urea	

Entry	Name	Structure
221	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-(trifluoromethyl)phenyl]carbamate	
224	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[5-(methylthio)pyridin-2-yl]phenyl}methyl)urea	
225	[3-(2,6-dimethylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
226	{3-[5-(methoxy)pyridin-3-yl]phenyl}methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
229	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-isoquinolin-4-yl)phenyl]methyl]urea	

Entry	Name	Structure
230	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-isoquinolin-4-yl)phenyl]methyl]urea	
232	[3-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
233	[4-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

28. (previously presented) A pharmaceutical composition comprising the compound according to claim 1 and a pharmaceutically acceptable carrier.

29. (cancelled)

30. (withdrawn from consideration, currently amended) A method for modulating the *in-vivo* activity of a kinase, the method comprising administering to a subject an effective amount of the compound according to ~~claim 1 to claim 1, or a compound selected from N-naphthalen-1-yl-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}-acetamide, N-[4-(phenyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}-acetamide, N-(3,4-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}~~

~~acetamide, N-(2,3-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,4-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,5-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(3,5-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,6-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-ethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,6-diethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2-(ethyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[3-(ethyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2,4-bis(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[4-(dimethylamino)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,3-dichlorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-chloro-3-methylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-bromophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2-fluorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-fluorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-N-[2-(trifluoromethyl)phenyl]-acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-N-[3-(trifluoromethyl)phenyl]-acetamide, methyl-4-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetyl]amino]-benzoate, ethyl-4-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetyl]amino]-benzoate, 3-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetyl]amino]-benzoic acid, N-[3-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[4-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy]-acetamide, N-(4-chlorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-aminophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, and N-(4-acetylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide.~~

31. (withdrawn from consideration) The method according to claim 30, wherein the kinase is c-Kit.

32. (withdrawn from consideration) The method according to claim 31, wherein modulating the *in vivo* activity of c-Kit comprises inhibition of c-Kit.

33. (withdrawn from consideration, currently amended) A method of treating rheumatoid arthritis, graft-host diseases, multiple sclerosis, psoriasis, arteriosclerosis, myocardioinfarction, ischemia, stroke, restenosis; interbowel diseases, osteoarthritis, macular degeneration, or diabetic retinopathy, ~~diseases or disorders associated with uncontrolled, abnormal, and/or unwanted cellular activities~~, the method comprising administering, to a mammal in need thereof, a therapeutically effective amount of the compound or the pharmaceutical composition as described in claim 1 ~~or a compound, or a pharmaceutical composition comprising said compound, selected from~~ N-naphthalen-1-yl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[4-(phenyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(3,4-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,3-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,4-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,5-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(3,5-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,6-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-N-(2,4,6-trimethylphenyl)-acetamide, N-(2-ethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-ethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,6-diethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2-(ethyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[3-(ethyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[2,4-bis(methyl-oxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-[4-(dimethylamino)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2,3-dichlorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-chloro-3-methylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-bromophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(2-fluorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, N-(4-fluorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-N-[2-(trifluoro-methyl)phenyl]-acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-N-[3-(trifluoromethyl)-phenyl]-acetamide, methyl-4-[[3-(1H-tetrazol-1-

yl)phenyl]oxy}acetyl)amino] — benzoate, — ethyl — 4-(((3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] — benzoate, 3-(((3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-[4-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-[2-chloro-5-(trifluoromethyl)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-(4-chlorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-(4-aminophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — and — N-(4-acetylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide.

34. (withdrawn from consideration, currently amended) A method of screening for modulators of c-Kit, the method comprising combining the compound according to claim 1 or a compound selected from N-naphthalen-1-yl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-[4-(phenyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-(3,4-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-(2,3-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-(2,4-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-(2,5-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-(3,5-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-(2,6-dimethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-(4-ethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-(2,6-diethylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-[2-(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-[2-(ethyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-[3-(ethyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-[2,4-bis(methyloxy)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-[4-(dimethylamino)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-(2,3-dichlorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-(4-chloro-3-methylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-(4-bromophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-(2-fluorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, — N-(4-fluorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, 2-[[3-(1H-

tetrazol-1-yl)phenyl]oxy} N-[2-(trifluoro-methyl)phenyl]—acetamide, —2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} N-[3-(trifluoromethyl)—phenyl]—acetamide, —methyl-4-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]—benzoate, —ethyl—4-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]—benzoate, —3-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, N-[4-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, —N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, —N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy}—acetamide, —N-(4-chlorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, N-(4-aminophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, —and—N-(4-acetylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, and at least one candidate agent and determining the effect of the candidate agent on c-Kit activity.

35. (withdrawn from consideration, currently amended) A method of inhibiting proliferative activity in a cell, the method comprising administering an effective amount of a composition comprising the compound according to claim 1 ~~or a compound selected from~~ N-naphthalen-1-yl-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, N-[4-(phenyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, —N-(3,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, —N-(2,3-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, —N-(2,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, N-(2,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, —N-(3,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, —N-(2,6-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, —2-{{[3-(1H-tetrazol-1-yl)phenyl]—oxy} N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, N-(4-ethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, N-(2,6-diethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, —N-[2-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, —N-[2-(ethyloxy)—phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, —N-[3-(ethyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, N-[4-(dimethylamino)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}—acetamide, —N-(2,3-

~~dichlorophenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(4-chloro-3-methylphenyl)-~~
~~2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(4-bromophenyl) 2-[[3-(1H-tetrazol-1-~~
~~yl)phenyl]oxy] acetamide, N-(2-fluorophenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide,~~
~~N-(4-fluorophenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, 2-[[3-(1H-tetrazol-1-~~
~~yl)phenyl]oxy] N-[2-(trifluoro-methyl)phenyl] acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]-~~
~~N-[3-(trifluoromethyl)-phenyl] acetamide, methyl 4-[[[3-(1H-tetrazol-1-~~
~~yl)phenyl]oxy]acetyl]amino] benzoate, ethyl 4-[[[3-(1H-tetrazol-1-~~
~~yl)phenyl]oxy]acetyl]amino] benzoate, 3-[[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetyl]amino]~~
~~benzoic acid, N-[3-(methyloxy)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-[4-~~
~~(methyloxy)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-[2-chloro-5-~~
~~(trifluoromethyl)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-[4-chloro-3-~~
~~(trifluoromethyl)phenyl] 2-[[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy] acetamide, N-(4-~~
~~chlorophenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(4-aminophenyl) 2-[[3-(1H-~~
~~tetrazol-1-yl)phenyl]oxy] acetamide, and N-(4-acetylphenyl) 2-[[3-(1H-tetrazol-1-~~
~~yl)phenyl]oxy] acetamide, to a cell or a plurality of cells.~~